

Data Path : Z:\voasrv\HPCHEM1\MSVOA_N\Data\VN040422\
 Data File : VN071835.D
 Acq On : 04 Apr 2022 16:41
 Operator : JC\MD
 Sample : N2175-01
 Misc : 5.0mL/MSVOA_N/WATER
 ALS Vial : 16 Sample Multiplier: 1

Instrument :
 MSVOA_N
 ClientSampleId :
 023-OUTFALL

Quant Time: Apr 05 01:19:18 2022
 Quant Method : Z:\voasrv\HPCHEM1\MSVOA_N\methods\82N033022W.M
 Quant Title : SW846 8260
 QLast Update : Thu Mar 31 01:57:54 2022
 Response via : Initial Calibration

Compound	R.T.	QIon	Response	Conc	Units	Dev(Min)

Internal Standards						
1) Pentafluorobenzene	8.080	168	637363	50.000	ug/l	0.00
34) 1,4-Difluorobenzene	8.963	114	1091083	50.000	ug/l	0.00
63) Chlorobenzene-d5	11.739	117	1113550	50.000	ug/l	0.00
72) 1,4-Dichlorobenzene-d4	13.668	152	423567	50.000	ug/l	0.00
System Monitoring Compounds						
33) 1,2-Dichloroethane-d4	8.433	65	359160	49.361	ug/l	0.00
Spiked Amount 50.000	Range 61 - 141		Recovery =	98.720%		
35) Dibromofluoromethane	8.016	113	310468	48.356	ug/l	0.00
Spiked Amount 50.000	Range 69 - 133		Recovery =	96.720%		
50) Toluene-d8	10.439	98	1376479	52.409	ug/l	0.00
Spiked Amount 50.000	Range 65 - 126		Recovery =	104.820%		
62) 4-Bromofluorobenzene	12.727	95	479226	56.417	ug/l	0.00
Spiked Amount 50.000	Range 58 - 135		Recovery =	112.840%		
Target Compounds						
					Qvalue	
13) Acrolein	4.051	56	4864	3.378	ug/l	90
16) Acetone	4.287	43	106913	34.382	ug/l	100
25) 2-Butanone	7.339	43	43081	9.610	ug/l	95
51) 4-Methyl-2-Pentanone	10.327	43	16578	1.767	ug/l	99
70) Styrene	12.292	104	36696	1.682	ug/l	89

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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